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                 CAOLD to be discontinued on December 31, 2008
NEWS 24
         AUG 15
                 CAplus currency for Korean patents enhanced
NEWS 25
        AUG 25
                 CA/CAplus, CASREACT, and IFI and USPAT databases
                 enhanced for more flexible patent number searching
NEWS 26
        AUG 27
                 CAS definition of basic patents expanded to ensure
                 comprehensive access to substance and sequence
                 information
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chain bonds :
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exact/norm bonds :
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exact bonds :
2-3 3-13
G1:C,O,N
G2:C,O,N
Match level :
1:CLASS 2:CLASS 3:CLASS 4:Atom 6:CLASS 7:CLASS 9:CLASS 10:Atom 12:Atom
13:CLASS 14:CLASS 15:CLASS
Element Count :
Node 10: Limited
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   N,N2
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L1 STRUCTURE UPLOADED

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SAMPLE SEARCH INITIATED 10:44:27 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 15428 TO ITERATE

13.0% PROCESSED 2000 ITERATIONS INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED) SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
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3 ANSWERS

PROJECTED ITERATIONS: 301120 TO 316000 PROJECTED ANSWERS: 174 TO 750

L2 3 SEA SSS SAM L1

=> d scan

L2 3 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN

IN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-[[2-(1-piperidinyl)phenyl]methyl]-

MF C25 H26 N6 O

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L2 3 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN

IN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-(5-methyl-3isoxazolyl)-

MF C17 H14 N6 O2

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L2 3 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN

IN Urea, N-[3-[2-(6-amino-3-pyridazinyl)ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]-

MF C20 H20 N6 O2

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> s c4n2/rf

L3 2537685 C4N2/RF

=> s 11 sub=13 full

FULL SUBSET SEARCH INITIATED 10:45:27 FILE 'REGISTRY'
FULL SUBSET SCREEN SEARCH COMPLETED - 26775 TO ITERATE

100.0% PROCESSED 26775 ITERATIONS 307 ANSWERS

SEARCH TIME: 00.00.01

L4 307 SEA SUB=L3 SSS FUL L1

=> file caplus

COST IN U.S. DOLLARS
SINCE FILE TOTAL
ENTRY SESSION
FULL ESTIMATED COST
183.97
184.18

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L5 8 L4

L5 ANSWER 8 OF 8 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 2005:158661 CAPLUS

DOCUMENT NUMBER: 142:240460

TITLE: Preparation of pyrimidine derivatives as ErbB kinase

inhibitors

INVENTOR(S): Reno, Michael John; Stevens, Kirk Lawrence; Waterson,

Alex Gregory; Zhang, Yuemei

PATENT ASSIGNEE(S): Smithkline Beecham Corporation, USA

SOURCE: PCT Int. Appl., 132 pp.

CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

GΙ

P.F	PATENT NO.				KIND DATE				APPLICATION NO.				DATE				
WC	WO 2005016914			A1 20050224			WO 2004-US26251				20040811						
	W:	ΑE,	AG,	AL,	AM,	ΑT,	ΑU,	AZ,	BA,	BB,	BG,	BR,	BW,	BY,	BZ,	CA,	CH,
		CN,	CO,	CR,	CU,	CZ,	DE,	DK,	DM ,	DZ,	EC,	EE,	EG,	ES,	FI,	GB,	GD,
		GE,	GH,	GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	KE,	KG,	KP,	KR,	KZ,	LC,
		LK,	LR,	LS,	LT,	LU,	LV,	MA,	MD,	MG,	MK,	MN,	MW,	MX,	MZ,	NA,	NI,
		NO,	NZ,	OM,	PG,	PH,	PL,	PT,	RO,	RU,	SC,	SD,	SE,	SG,	SK,	SL,	SY,
		ТJ,	TM,	TN,	TR,	TT,	TZ,	UA,	UG,	US,	UΖ,	VC,	VN,	YU,	ZA,	ZM,	ZW
	RW:	BW,	GH,	GM,	ΚE,	LS,	MW,	MZ,	NA,	SD,	SL,	SZ,	TZ,	UG,	ZM,	ZW,	AM,
		AZ,	BY,	KG,	KZ,	MD,	RU,	ТJ,	TM,	AT,	BE,	BG,	CH,	CY,	CZ,	DE,	DK,
		EE,	ES,	FI,	FR,	GB,	GR,	HU,	ΙE,	IT,	LU,	MC,	NL,	PL,	PT,	RO,	SE,
		SI,	SK,	TR,	BF,	ВJ,	CF,	CG,	CI,	CM,	GA,	GN,	GQ,	GW,	ML,	MR,	NE,
		SN,	TD,	TG													
EF	EP 1654251			A1 20060510				EP 2004-781004				20040811					
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JE	JP 2007502298				T 20070208 JP 2006-523388						88	20040811					
US	US 20060205740			A1 20060914				US 2006-568052				20060210					
PRIORIT	PRIORITY APPLN. INFO.:				US 2003-495180P						I	P 20030814					
									,	WO 2	004-1	US26	251	Ţ	W 2	0040	811
OTHER SOURCE(S): CF					CASI	CASREACT 142:240460; MARPAT 142:240460											

Title compds. I [wherein A = alkenylene, alkynylene; R = alkylene; R1 = AΒ -(Z)-(Z1)m-(Z2)n; Z = hetero/aryl, hetero/arylene; Z1 = CH2 where m = 0-1; Z2 = OH and derivs., halo, CN, CONH2 and derivs. or heterocyclyl, where n = 0-1, etc.; R2 = H, alkyl; R3 = -(Q)-(Q1)r-(Q2); Q = hetero/arylene; Q1 = O, where r = 0-1; Q2 = arylalkyl, hetero/aryl; and their salts, solvates, and physiol. functional derivs.] were prepared as ErbB kinase inhibitors for treating cancer. Thus, reacting 2-benzyl-N-(5-vinylpyrimidin-4-yl)-1Hbenzimidazol-5-amine (preparation given) with Ph iodide gave pyrimidine II in 8%. I showed inhibitory activity vs. EGFR, ErbB-2, and ErbB-4 protein tyrosine kinases with a pIC50 \geq 5.0. I are useful in the treatment of diseases associated with inappropriate ErbB family kinase activity. 845657-86-5P, N-[3-Chloro-4-[(3-fluorobenzyl)oxy]phenyl]-5-[[6-[[[2-(4-morpholinyl)ethyl]amino]methyl]-2-pyridinyl]ethynyl]-4pyrimidinamine 845657-90-1P, N-[3-Chloro-4-[(3fluorobenzyl)oxy]phenyl]-5-[[6-[[[3-(1H-imidazol-1-yl)propyl]amino]methyl]-2-pyridinyl]ethynyl]-4-pyrimidinamine 845658-10-8P, 1-[[6-[[4-[3-Chloro-4-[(3-fluorobenzyl)oxy]anilino]-5-pyrimidinyl]ethynyl]-2-pyridinyl]methyl]-3-[2-(4-morpholinyl)ethyl]urea RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses) (drug candidate; preparation of pyrimidines as ErB kinase inhibitors) RN 845657-86-5 CAPLUS 4-Morpholineethanamine, N-[[6-[2-[4-[[3-chloro-4-[(3-CN fluorophenyl)methoxy]phenyl]amino]-5-pyrimidinyl]ethynyl]-2pyridinyl]methyl]- (CA INDEX NAME)

ΙI

RN 845657-90-1 CAPLUS

CN 4-Pyrimidinamine, N-[3-chloro-4-[(3-fluorophenyl)methoxy]phenyl]-5-[2-[6-[[3-(1H-imidazol-1-yl)propyl]amino]methyl]-2-pyridinyl]ethynyl]- (CA INDEX NAME)

RN 845658-10-8 CAPLUS

CN Urea, N-[[6-[2-[4-[[3-chloro-4-[(3-fluorophenyl)methoxy]phenyl]amino]-5-pyrimidinyl]ethynyl]-2-pyridinyl]methyl]-N'-[2-(4-morpholinyl)ethyl]- (CA INDEX NAME)

PAGE 2-A



REFERENCE COUNT: 1 THERE ARE 1 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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L5 ANSWER 7 OF 8 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 2005:588667 CAPLUS

DOCUMENT NUMBER: 143:115556

TITLE: Preparation of 4-aminopyrimidine derivatives as

inhibitors of Tie2 receptor tyrosine kinases

INVENTOR(S): Jones, Clifford David; Luke, Richard William Arthur;

McCoull, William

PATENT ASSIGNEE(S): Astrazeneca AB, Swed.; Astrazeneca UK Limited

SOURCE: PCT Int. Appl., 129 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2005060969	A1	20050707	WO 2004-GB5332	20041220

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             LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI,
             NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY,
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     EP 1737462
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     IN 2006MN00847
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PRIORITY APPLN. INFO.:
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                                                                     20031224
                                             GB 2004-16850
                                                                  Α
                                                                     20040729
                                             WO 2004-GB5332
                                                                  W
                                                                     20041220
OTHER SOURCE(S):
                         CASREACT 143:115556; MARPAT 143:115556
GΙ
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AB Title compds. I [wherein R1, R2 = H, alkyl, alkanoyl; R3, R4 = H, alkyl, alkoxy; R5 = cyclopropyl, halo, cyano; m, n = 0-3; R6 = halo, oxo, cyano; etc., or salts thereof] were prepared as inhibitors of Tie2 receptor tyrosine kinases. Processes for the synthesis of I and some intermediates involved are claimed. For example, urea II was synthesized in 21% yield by condensation of the corresponding aniline with Ph thiadiazolylcarbamate in the presence of Et3N in THF under microwave irradiation. This urea showed inhibition against Tie2 receptor tyrosine kinase in vitro and inhibition of autophosphorylation of Tie2 receptor tyrosine kinase with IC50 values of 0.879 μM and 5.557 μM, resp. Therefore, I and their pharmaceutical compns. have potential use in the production of an anti-angiogenic effect in a warm-blooded animal.

IT 857287-13-9P, Phenyl [3-[(4,6-diaminopyrimidin-5-

RN 857287-53-7 CAPLUS

CN Urea, N-[3-(1,1-dimethylethyl)-1-methyl-1H-pyrazol-5-yl]-N'-[3-[2-[4-[[2-(4-morpholinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)

ΙT 857287-02-6P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'phenylurea 857287-04-8P, 2-Phenyl-N-[3-[(4,6-diaminopyrimidin-5yl)ethynyl]phenyl]acetamide 857287-05-9P, N-[3-[(4,6- $\label{lem:diaminopyrimidin-5-yl} Diaminopyrimidin-5-yl) ethynyl] phenyl]-N'-(3,4-dichlorophenyl) urea$ 857287-06-0P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-[2-(trifluoromethyl)phenyl]urea 857287-07-1P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-[3-(trifluoromethyl)phenyl]urea 857287-08-2P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-[4-(trifluoromethyl)phenyl]urea [2-fluoro-5-(trifluoromethyl)phenyl]urea 857287-10-6P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(3-methoxyphenyl)urea 857287-11-7P, Phenyl [4-[(4,6-diaminopyrimidin-5v1) ethynyl]phenyl]carbamate 857287-14-0P, N-(5-tert-Butyl-1,3,4thiadiazol-2-yl)-N'-[4-[(4,6-diaminopyrimidin-5-yl)ethynyl]phenyl]urea 857287-15-1P, N-[4-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(3-methylisothiazol-5-yl)urea 857287-16-2P, N-[4-[(4,6-4)]]Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(3-methylisoxazol-5-yl)urea 857287-17-3P, N-[4-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-[4-(trifluoromethyl)pyridin-2-yl]urea 857287-18-4P, N-[3-[[[4-[4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]amino]carbonyl]amino]phenyl]acetamide 857287-19-5P, N-[3-[(4,6-Diaminopyrimidin-5y1)ethynyl]phenyl]-N'-(3-methylisothiazol-5-y1)urea 857287-20-8P , N-[3-[[[[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]amino]carbonyl]ami

```
no]phenyl]acetamide 857287-21-9P, N-[3-[(4,6-Diaminopyrimidin-5-
yl)ethynyl]phenyl]-N'-[4-(trifluoromethyl)pyridin-2-yl]urea
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857287-24-2P, N-(5-tert-Butylisoxazol-3-y1)-N'-[3-[(4,6-
diaminopyrimidin-5-yl)ethynyl]phenyl]urea 857287-25-3P,
N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(2,3-dihydro-1,4-
benzodioxin-6-yl)urea 857287-26-4P, N-[3-[(4,6-Diaminopyrimidin-
5-yl)ethynyl]phenyl]-N'-[2-(morpholin-4-yl)phenyl]urea
857287-27-5P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-
(1-methylpiperidin-4-y1)urea 857287-28-6P, N-[3-[(4,6-
Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(1-propylpiperidin-4-yl)urea
857287-29-7P, N-[3-[(4,6-Diaminopyrimidin-5-y1)ethynyl]phenyl]-2-
(2-methoxypheny1) acetamide 857287-30-0P, N-[3-[(4,6-
Diaminopyrimidin-5-yl)ethynyl]phenyl]-2-[3-(trifluoromethyl)phenyl]acetami
de 857287-31-1P, N-[3-[(4,6-Diaminopyrimidin-5-
yl)ethynyl]phenyl]-2-[4-(trifluoromethyl)phenyl]acetamide
857287-32-2P, N-[3-[(4,6-Diaminopyrimidin-5-y1)ethynyl]phenyl]-2-
(3-methoxyphenyl)acetamide 857287-35-5P, N-(5-tert-Butylisoxazol-
3-y1)-N'-[3-[[4-(methylamino)pyrimidin-5-y1]ethynyl]phenyl]urea
857287-36-6P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[4-[[3-yl]]]]
(isopropylamino)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea
857287-37-7P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[4-[[2-1]]]]
(pyrrolidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea
857287-38-8P, N-(5-tert-Butylisoxazol-3-y1)-N'-[3-[[4-[(5-tert-
butylisoxazol-3-yl)amino]pyrimidin-5-yl]ethynyl]phenyl]urea
(dimethylamino)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea
857287-40-2P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[4-[(2-
hydroxyethyl)amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857287-41-3P
, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[4-[[2-(morpholin-4-
y1)ethy1]amino]pyrimidin-5-y1]ethyny1]pheny1]urea 857287-42-4P,
N-[3-[[4-[(4-Aminobutyl)amino]pyrimidin-5-yl]ethynyl]phenyl]-N'-(5-tert-
butylisoxazol-3-yl)urea 857287-43-5P, N-(5-tert-Butylisoxazol-3-
y1)-N'-[3-[4-[3-(pyrrolidin-1-y1)propy1]amino]pyrimidin-5-
yl]ethynyl]phenyl]urea 857287-44-6P, N-(5-tert-Butylisoxazol-3-
y1)-N'-[3-[4-(2,4-dimethoxybenzyl)amino]pyrimidin-5-
yl]ethynyl]phenyl]urea 857287-45-7P, N-[3-[[4-[(2-
Aminoethyl)amino]pyrimidin-5-yl]ethynyl]phenyl]-N'-(5-tert-butylisoxazol-3-
y1)urea 857287-46-8P, N-(5-tert-Butylisoxazol-3-y1)-N'-[3-[[4-
[[2-(dimethylamino)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea
(dimethylamino)butyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea
857287-48-0P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[4-[N-[2-
(dimethylamino)ethyl]methylamino]pyrimidin-5-yl]ethynyl]phenyl]urea
857287-49-1P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[4-[[2-1]]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]]-N'-[3-[[4-1]]
(piperidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea
857287-50-4P, N-(5-tert-Butylisoxazol-3-y1)-N'-[3-[[4-[[3-
(morpholin-4-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea
857287-51-5P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[4-[[3-
(piperidin-1-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea
857287-52-6P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[4-[[3-(4-1)]]]]
methylpiperazin-1-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea
857287-57-1P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-
(2,3-dihydro-1H-inden-1-yl)urea
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU
(Therapeutic use); BIOL (Biological study); PREP (Preparation); USES
(Uses)
      (inhibitor; preparation of aminopyrimidine derivs. as inhibitors of Tie2
      receptor tyrosine kinases)
```

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-phenyl- (CA INDEX NAME)

RN 857287-04-8 CAPLUS

CN Benzeneacetamide, N-[3-[2-(4,6-diamino-5-pyrimidiny1)ethynyl]phenyl]- (CA INDEX NAME)

$$C = C$$
 $NH - C - CH_2 - Ph$
 $NH - C - CH_2 - Ph$
 $NH - C - CH_2 - Ph$

RN 857287-05-9 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(3,4-dichlorophenyl)- (CA INDEX NAME)

RN 857287-06-0 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-[2-(trifluoromethyl)phenyl]- (CA INDEX NAME)

RN 857287-07-1 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-[3-(trifluoromethyl)phenyl]- (CA INDEX NAME)

RN 857287-08-2 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-[4-(trifluoromethyl)phenyl]- (CA INDEX NAME)

RN 857287-09-3 CAPLUS

CN Urea, N-[4-[2-(4,6-diamino-5-pyrimidiny1)ethyny1]pheny1]-N'-[2-fluoro-5-(trifluoromethy1)pheny1]- (CA INDEX NAME)

RN 857287-10-6 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(3-methoxyphenyl)- (CA INDEX NAME)

$$N$$
 NH_2
 NH_2
 NH_2
 NH_2
 NH_2
 NH_2
 NH_2

RN 857287-11-7 CAPLUS

CN Carbamic acid, [4-[(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-, phenyl ester (9CI) (CA INDEX NAME)

$$NH_2$$
 NH_2
 NH_2
 NH_2
 NH_2
 NH_2

RN 857287-14-0 CAPLUS

CN Urea, N-[4-[2-(4,6-diamino-5-pyrimidiny1)ethyny1]pheny1]-N'-[5-(1,1-dimethylethyl)-1,3,4-thiadiazol-2-yl]- (CA INDEX NAME)

RN 857287-15-1 CAPLUS

CN Urea, N-[4-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(3-methyl-5-isothiazolyl)- (CA INDEX NAME)

RN 857287-16-2 CAPLUS

CN Urea, N-[4-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(3-methyl-5-isoxazolyl)- (CA INDEX NAME)

RN 857287-17-3 CAPLUS

CN Urea, N-[4-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-[4-(trifluoromethyl)-2-pyridinyl]- (CA INDEX NAME)

RN 857287-18-4 CAPLUS

CN Acetamide, N-[3-[[[[4-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]amino]c arbonyl]amino]phenyl]- (CA INDEX NAME)

RN 857287-19-5 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(3-methyl-5-isothiazolyl)- (CA INDEX NAME)

RN 857287-20-8 CAPLUS

CN Acetamide, N-[3-[[[[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]amino]c arbonyl]amino]phenyl]- (CA INDEX NAME)

RN 857287-21-9 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-[4-(trifluoromethyl)-2-pyridinyl]- (CA INDEX NAME)

RN 857287-22-0 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(3-methyl-5-isoxazolyl)- (CA INDEX NAME)

RN 857287-23-1 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-1,3,4-thiadiazol-2-yl]- (CA INDEX NAME)

RN 857287-24-2 CAPLUS

CN Urea, $N-[3-[2-(4,6-diamino-5-pyrimidinyl)] = 10^{-1} - 10^{-1$

RN 857287-25-3 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(2,3-dihydro-1,4-benzodioxin-6-yl)- (CA INDEX NAME)

RN 857287-26-4 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidiny1)ethyny1]pheny1]-N'-[2-(4-morpholiny1)pheny1]- (CA INDEX NAME)

RN 857287-27-5 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(1-methyl-4-piperidinyl)- (CA INDEX NAME)

$$\begin{array}{c|c} NH_2 & O & Me \\ \hline N & NH_2 & NH-C-NH & NH_2 & NH_$$

RN 857287-28-6 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(1-propyl-4-piperidinyl)- (CA INDEX NAME)

RN 857287-29-7 CAPLUS

CN Benzeneacetamide, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-2-methoxy- (CA INDEX NAME)

$$\begin{array}{c|c} N & NH_2 & O & MeO \\ \hline N & C & \hline \end{array} \\ NH_2 & NH_2 & O & MeO \\ \hline \end{array}$$

RN 857287-30-0 CAPLUS

CN Benzeneacetamide, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-3-(trifluoromethyl)- (CA INDEX NAME)

$$\begin{array}{c|c} N & NH_2 & O \\ N & C \longrightarrow C \end{array}$$

$$NH_2 & C \longrightarrow C \longrightarrow C$$

$$NH_2 & C \longrightarrow C$$

$$CF_3$$

RN 857287-31-1 CAPLUS

CN Benzeneacetamide, N-[3-[2-(4,6-diamino-5-pyrimidiny1)]] ethynyl]phenyl]-4-(trifluoromethyl)- (CA INDEX NAME)

$$\begin{array}{c|c} NH_2 & O \\ NH_2 & C \\ \hline NH_2 & NH_2 & C \\ \end{array}$$

RN 857287-32-2 CAPLUS

CN Benzeneacetamide, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-3-methoxy- (CA INDEX NAME)

RN 857287-35-5 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[4-(methylamino)-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)

RN 857287-36-6 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[4-[[3-[(1-methylethyl)amino]propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)

RN 857287-37-7 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[4-[[2-(1-pyrrolidinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)

RN 857287-38-8 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[4-[[5-(1,1-dimethylethyl)-3-isoxazolyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)

$$\begin{array}{c} Bu-t \\ N \\ N \\ C \end{array} \begin{array}{c} C \\ NH-C-NH \\ O \\ Bu-t \end{array}$$

ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

RN 857287-39-9 CAPLUS

CN Urea, N-[3-[2-[4-[[3-(dimethylamino)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)

$$\begin{array}{c} & \text{Me}_2N - (CH_2)_3 - NH \\ & \text{NH-C-NH} \\ & \text{t-Bu} \end{array}$$

RN 857287-40-2 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[4-[(2-iyana)]]-N'-[3-[2-[4-[(2-iyana)]]]-N'-[3-[2-[4-[(2-[4-[(2-[4-[iyana)]]]]]-N'-[3-[2-[4-[(2-[4-[iyana)]]]]]-N'-[3-[4-[iyana]]]-N'-[3-[4-[iyana]]]-N'-[4-[iyana]]-N'-[4-[i

hydroxyethyl)amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)

RN 857287-41-3 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[4-[[2-(4-morpholinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)

RN 857287-42-4 CAPLUS

CN Urea, N-[3-[2-[4-[(4-aminobutyl)amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)

RN 857287-43-5 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[4-[[3-(1-pyrrolidinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)

RN 857287-44-6 CAPLUS

CN Urea, N-[3-[2-[4-[[(2,4-dimethoxyphenyl)methyl]amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)

RN 857287-45-7 CAPLUS

CN Urea, N-[3-[2-[4-[(2-aminoethyl)amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)

RN 857287-46-8 CAPLUS

CN Urea, N-[3-[2-[4-[[2-(dimethylamino)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)

RN 857287-47-9 CAPLUS

CN Urea, N-[3-[2-[4-[[4-(dimethylamino)butyl]amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)

RN 857287-48-0 CAPLUS

CN Urea, N-[3-[2-[4-[[2-(dimethylamino)ethyl]methylamino]-5-pyrimidinyl]ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)

t-Bu
$$Me_2N-CH_2-CH_2-N$$

RN 857287-49-1 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[4-[[2-(1-piperidinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)

RN 857287-50-4 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[4-[[3-(4-morpholinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)

RN 857287-51-5 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[4-[[3-(1-piperidinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)

RN 857287-52-6 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[4-[[3-(4-methyl-1-piperazinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)

Me N
$$\sim$$
 (CH2)3-NH \sim N \sim C \sim C \sim N \sim NH-C-NH

RN 857287-57-1 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidiny1)ethyny1]pheny1]-N'-(2,3-dihydro-1H-inden-1-y1)- (CA INDEX NAME)

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